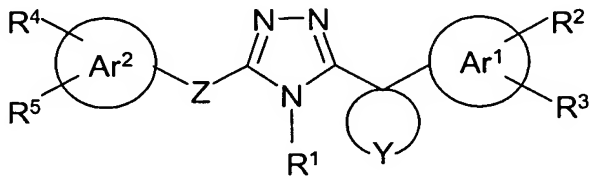
This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. **(Original)** A triazole compound represented by the following formula:



wherein

R^1 is an alkyl group or a cycloalkyl group

wherein the alkyl group and the cycloalkyl group are optionally substituted by 1 to 5 substituents each independently selected from a halogen atom, $-CF_3$, -OH,

$-NH_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, -COOH, -CO-O-alkyl, $-CO-N(R^7)(R^8)$, $-N(R^7)-CO-R^8$, an aryl group and a heteroaryl group

wherein R^7 and R^8 are each independently a hydrogen atom or an alkyl group, and the aryl group and the heteroaryl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group,

$-(CH_2)_n-OH$, $-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group

wherein n is 0-3, R^9 and R^{10} are each independently a hydrogen atom, an alkyl group or –CO-alkyl, and R^{11} is -OH, an alkoxy group, an alkyl group or

$-N(R^{12})(R^{13})$ wherein R^{12} and R^{13} are each independently a hydrogen atom or an alkyl group;

Y is a cycloalkyl group or a heterocycloalkyl group

wherein the cycloalkyl group and the heterocycloalkyl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group,

$-(CH_2)_n-OH$, $-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above);

Ar^1 is an aryl group or a heteroaryl group;

R^2 and R^3

are each independently a hydrogen atom, a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n-OH$, $-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group or a heteroaryl group

wherein the aryl group and the heteroaryl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n-OH$,

$-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above);

Z is $-(CH(R^{14}))_p-$, $-(CH(R^{14}))_p-N(R^{16})-(CH(R^{15}))_q-$ or

wherein Y_1 is a cycloalkyl group or a heterocycloalkyl group

wherein the cycloalkyl group and the heterocycloalkyl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n-OH$, $-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above),

p is 0-3, q is 0-3, R^{14} and R^{15} are each independently a hydrogen atom, a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n-OH$, $-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group,

$-CO-R^{11}$, an aryl group or a heteroaryl group

wherein the aryl group and the heteroaryl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n-OH$,

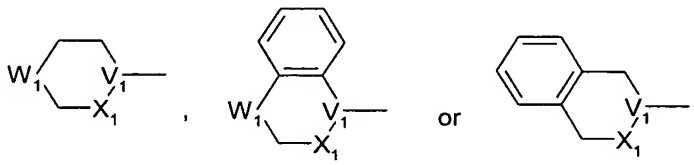
$-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above), and

R^{16} is a hydrogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n-OH$, $-(CH_2)_n-CO-R^{11}$, a cycloalkyl group, an alkenyl group, an aryl group or a heteroaryl group

wherein the aryl group and the heteroaryl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n-OH$,

$-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above);

Ar^2 is an aryl group or a heteroaryl group; or



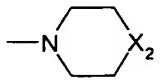
wherein X_1 is $-(CH_2)_t-$ wherein t is 0-2, V_1 is =CH- or =N-, and W_1 is $-C(R^{17})(R^{18})-$, -O-, -S-, $-SO_2-$, -SO-, -CO- or

$-N(R^{19})-$

wherein R^{17} and R^{18} are each independently a hydrogen atom, an alkyl group, an alkoxy group, a haloalkyl group, $-(CH_2)_r-OH$, $-CO-R^{20}$, $-N(R^{21})(R^{22})$ or $-L_1-Ar^3$

wherein r is 0-3, R^{20} is -OH, an alkoxy group, an alkoxyalkyl group or $-N(R^{23})(R^{24})$

wherein R^{23} and R^{24} are each independently a hydrogen atom, an alkyl group, $-(CH_2)_s$-OH, an alkoxyalkyl group, or in combination form



wherein s is 0-3, X_2 is -O-, $-(CH_2)_t$- or $-N(R^{25})$-

wherein t is as defined above and R^{25} is a hydrogen atom, $-CO-R^{26}$, $-SO_2-R^{26}$ or $-(CH_2)_u-Ar^4$

wherein R^{26} is an alkyl group, an alkoxy group, -NH-alkyl or $-N(\text{-alkyl})_2$, u is 0-3, and Ar^4 is an aryl group or a heteroaryl group wherein the aryl group and the heteroaryl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n$-OH,

$-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above),

L_1 is $-(CH_2)_v$-, -O- or -CO-

wherein v is 0-3, and

Ar^3 is an aryl group or a heteroaryl group wherein the aryl group and the heteroaryl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n$-OH, $-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group,

$-CO-R^{11}$, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above), and

R^{21} and R^{22} are each independently a hydrogen atom, an alkyl group, -CO-alkyl, -CO-O-alkyl or $-L_1-Ar^3$ (L_1 and Ar^3 are as defined above), and

R^{19} is a hydrogen atom, $-CO-R^{26}$, $-SO_2-R^{26}$ or $-(CH_2)_u-Ar^4$ (R^{26}, u and Ar^4 are as defined above); and

R^4 and R^5

are each independently a hydrogen atom, a halogen atom, -OH, $-NO_2$, -CN, an alkyl group, an alkoxy group, $-CO-R^{27}$, $-SO_2-R^{27}$, $-CO-N(R^{28})(R^{29})$ or $-N(R^{30})(R^{31})$

wherein the alkyl group and the alkoxy group are optionally substituted by 1 to 5 substituents each independently selected from a halogen atom, $-CF_3$, -OH, an alkoxy group, a haloalkoxy group, $-N(R^9)(R^{10})$, -CN, $-NO_2$, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group (R^9, R^{10} and R^{11} are as defined above),

wherein the aryl group and the heteroaryl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group, $-(CH_2)_n-OH$,

$-N(R^9)(R^{10})$, -CN, $-NO_2$, an alkoxy group, a cycloalkyl group, an alkenyl group, $-CO-R^{11}$, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above)

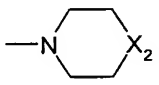
R^{27} is -OH, an alkoxy group, an alkyl group, $-NH_2$,

-NH-alkyl or $-N(-alkyl)_2$,

R^{28} and R^{29} are each independently a hydrogen atom, an alkyl group or $-(CH_2)_w-R^{32}$,

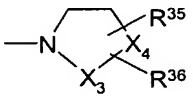
wherein w is 0-3 and R^{32} is -OH, $-CF_3$, an alkoxy group, $-CONH_2$ or $-N(R^{33})(R^{34})$

wherein R^{33} and R^{34} are each independently a hydrogen atom, an alkyl group, -CO-alkyl, or in combination form



(X_2 is as defined above)

or R^{28} and R^{29} in combination form

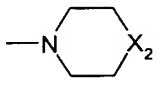


wherein X_3 is -CO-, $-CH_2-$ or $-CH_2-CH_2-$, X_4 is

-O-, -$(CH_2)_t$-, -$N(R^{25})$- or 

wherein Y_2 is cycloalkyl or heterocycloalkyl and t and R^{25} are as defined above, and R^{35} and R^{36} are each independently a hydrogen atom, a halogen atom, an alkyl group optionally substituted by -OH, -OH, -CN, -NO_2, an alkoxy group, a cycloalkyl group, an alkenyl group, -CO-R^{37}, -$N(R^{38})(R^{39})$

wherein R^{37} is -OH, an alkoxy group, -NH_2,

-NH-alkyl, -$N(\text{-alkyl})_2$ or  (X_2 is as defined above)

wherein the alkyl group in –NH-alkyl and -$N(\text{-alkyl})_2$ and the alkoxy group are optionally substituted by 1 to 5 substituents each independently selected from a halogen atom, -CF_3, -OH, an alkoxy group, a haloalkoxy group,

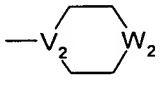
-$N(R^9)(R^{10})$, -CN, -NO_2, a cycloalkyl group, an alkenyl group,

-CO-R^{11}, an aryl group and a heteroaryl group (R^9, R^{10} and R^{11} are as defined above),

wherein the aryl group and the heteroaryl group are optionally substituted by 1 to 3 substituents each independently selected from a halogen atom, a haloalkyl group, an alkyl group, -$(CH_2)_n$-OH, -$N(R^9)(R^{10})$, -CN, -NO_2, an alkoxy group, a cycloalkyl group, an alkenyl group, -CO-R^{11}, an aryl group and a heteroaryl group (n, R^9, R^{10} and R^{11} are as defined above), and

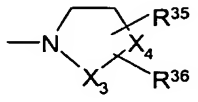
R^{38} and R^{39} are each independently a hydrogen atom, an alkyl group, -CO-alkyl or -CO-O-alkyl, and

R^{30} and R^{31} are each independently a hydrogen atom, an alkyl group optionally substituted by

–OH, -SO_2-R^{40}, -$(CH_2)_x$-CO-R^{41} or 

wherein x is 0-3, R^{40} is an alkyl group or -NH_2, R^{41} is a hydrogen atom, an alkyl group optionally substituted by -OH, -OH, an alkoxy group, an alkoxyalkyl group or -$(CH_2)_s$-$N(R^{42})(R^{43})$

wherein s is as defined above and R^{42} and R^{43} are each independently a hydrogen atom, an alkyl group, -OH, an alkoxy group, or in combination form


(X_3, X_4, R^{35} and R^{36} are as defined above),

V_2 is =CH- or =N- and W_2 is $-C(R^{44})(R^{45})-$, -O- or $-N(R^{46})-$

wherein R^{44} and R^{45} are each independently a hydrogen atom, an alkyl group, an alkoxy group, a haloalkyl group, $-(CH_2)_r-OH$, $-CO-R^{47}$ or $-N(R^{48})(R^{49})$

wherein r is as defined above, R^{47} is -OH, an alkoxy group, an alkoxyalkyl group,

$-N(R^{50})(R^{51})$

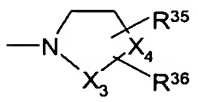
wherein R^{50} and R^{51} are each independently a hydrogen atom, an alkyl group, $-(CH_2)_s-OH$ (s is as defined above) or an alkoxyalkyl group, and

R^{48} and R^{49} are each independently a hydrogen atom, an alkyl group, -CO-alkyl or -CO-O-alkyl, and

R^{46} is a hydrogen atom, $-CO-R^{52}$ or $-SO_2-R^{52}$

wherein R^{52} is an alkyl group, an alkoxy group, -NH-alkyl or $-N(-alkyl)_2$ or

R^{30} and R^{31} in combination form


(X_3, X_4, R^{35} and R^{36} are as defined above), or

R^4 and R^5 in combination may form –O-alkylene-O-,

a prodrug thereof or a pharmaceutically acceptable salt thereof.

2. (Original) The triazole compound of claim 1, wherein Z is $-(CH(R^{14}))_p-$ and p is 0, a prodrug thereof or a pharmaceutically acceptable salt thereof.

3. (Original) The triazole compound of claim 2, wherein Y is a C_{3-8} cycloalkyl group, a prodrug thereof or a pharmaceutically acceptable salt thereof.

4. (Original) The triazole compound of claim 3, wherein Ar^1 is a phenyl group, a prodrug thereof or a pharmaceutically acceptable salt thereof.

5. (Original) The triazole compound of claim 4, wherein R^2 and R^3 are each independently a halogen atom or a hydrogen atom, a prodrug thereof or a pharmaceutically acceptable salt thereof.

6. (Previously Presented) The triazole compound of claim 1, wherein Ar^2 is a phenyl group, R^4 is a hydrogen atom and R^5 is $-CO-N(R^{28})(R^{29})$, a prodrug thereof or a pharmaceutically acceptable salt thereof.

7. (Original) The triazole compound of claim 6, wherein R^{28} and R^{29} are each independently a hydrogen atom or an alkyl group, a prodrug thereof or a pharmaceutically acceptable salt thereof.

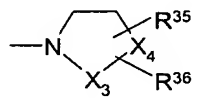
8. (Previously Presented) The triazole compound of claim 1, wherein Ar^2 is a phenyl group, R^4 is a hydrogen atom and R^5 is $-N(R^{30})(R^{31})$ wherein R^{30} is a hydrogen atom and R^{31} is $-(CH_2)_x-CO-R^{41}$, a prodrug thereof or a pharmaceutically acceptable salt thereof.

9. (Original) The triazole compound of claim 8, wherein X is 0 and R^{41} is an alkoxy group, a prodrug thereof or a pharmaceutically acceptable salt thereof.

10. (Original) The triazole compound of claim 8, wherein X is 0 and R^{41} is $-(CH_2)_s-N(R^{42})(R^{43})$, a prodrug thereof or a pharmaceutically acceptable salt thereof.

11. (Original) The triazole compound of claim 10, wherein s is 0, R^{42} is a hydrogen atom and R^{43} is an alkoxy group, a prodrug thereof or a pharmaceutically acceptable salt thereof.

12. (Previously Presented) The triazole compound of claim 1, wherein Ar^2 is a phenyl group, R^4 is a hydrogen atom and R^5 is $-N(R^{30})(R^{31})$ wherein R^{30} and R^{31} are joined to form



and X_3 is –CO-, a prodrug thereof or a pharmaceutically acceptable salt thereof.

13. (Original) The triazole compound of claim 12, wherein X_4 is –O-, a prodrug thereof or a pharmaceutically acceptable salt thereof.

14. (Original) The triazole compound of claim 1, which is

3-chloro-4-[4-methyl-5-(1-phenyl-cyclopropyl)-4H-[1,2,4]triazol-3-yl]-benzamide,

{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzoyl}morpholine hydrochloride,

3-chloro-N-methyl-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide hydrochloride,

3-chloro-N,N-dimethyl-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide hydrochloride,

3-chloro-N-(2-hydroxy-ethyl)-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-N-isopropyl-4-[4-methyl-5-(1-phenylcyclopropyl)4H-[1,2,4]triazol-3-yl]benzamide,

{3-chloro-4-[4-methyl-5-(1-phenyl-cyclopropyl)-4H[1,2,4]triazol-3-yl]benzoyl}piperidine hydrochloride,

{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H[1,2,4]triazol-3-yl]benzoyl}-(4-hydroxy)piperidine,

N-carbamoylmethyl-3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-N-(2,2,2-trifluoro-ethyl)-benzamide hydrochloride,

N-(2-acetylamino)ethyl-3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide hydrochloride,

3-chloro-N-(2-methoxy)ethyl-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide hydrochloride,

1-acetyl-(4-{3-Chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzoyl}piperazine hydrochloride,

3-chloro-N-(2-dimethylamino)ethyl-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-N-(2-morpholin-4-yl)ethylbenzamide,

4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-3-methoxybenzamide,

3-chloro-4-{4-methyl-5-[1-(4-fluorophenyl)cyclopropyl]-4H-[1,2,4]triazol-3-yl}benzamide,

3-chloro-N-metyl-4-{4-methyl-5-[1-(4-fluoro-phenyl)cyclopropyl]-4H-[1,2,4]triazol-3-yl}benzamide,

4-[4-isopropyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-4-[4-methyl-5-(1-thiophen-2-yl)cyclopropyl-4H-[1,2,4]triazol-3-yl]benzamide,

4-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}benzamide,

4-chloro-3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}benzamide,

4-chloro-3-{5-[1-phenylcyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}benzamide,

3-chloro-4-[4-ethyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-4-{4-ethyl-5-[1-(4-fluorophenyl)cyclopropyl]-4H-[1,2,4]triazol-3-yl}benzamide,

3-[4-isopropyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-{5-[1-(4-fluoro-phenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}benzamide,

N-{3-chloro-4-[4-methyl-5-(1-phenylcycloproppyl)-4H-[1,2,4]triazol-3-yl]phenyl}-1-morpholinecarboxamide,

3-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}-1,1-dimethylurea,

{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}urea,

ethyl N-{3-Chloro-4-[4-methyl-5-(1-phenyl-cyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}-carbamate,

N- { 3-chloro-4-[4-methyl-5-(1-phenylcycloproppyl)-4H-[1,2,4]triazol-3-yl]phenyl } -1-(4-methoxypiperidine)carboxamide,

N- { 3-chloro-4-[4-methyl-5-(1-phenylcycloproppyl)-4H-[1,2,4]triazol-3-yl]phenyl } -1-(3-hydroxypiperidine)carboxamide,

N- { 3-chloro-4-[4-methyl-5-(1-phenylcycloproppyl)-4H-[1,2,4]triazol-3-yl]phenyl } -1-(4-hydroxypiperidine)carboxamide,

1-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}-3-methoxyurea,

1-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]phenyl}-3-hydroxy-3-methylurea,

1-(3-chloro-4-{5-[1-(4-fluorophenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}phenyl)-3-methoxyurea,

1-(4-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}phenyl)-3-methoxyurea,

1-(3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}phenyl)-3-methoxyurea,

3-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}oxazolidin-2-one,

1-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]phenyl}imidazolidin-2-one,

3-(3-chloro-4-{5-[1-(4-fluoro-phenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}phenyl)oxazolidin-2-one,

3-(4-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}phenyl)oxazolidin-2-one,

3-(4-chloro-3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}phenyl)oxazolidin-2-one,

3-(3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}phenyl)oxazolidin-2-one,

methyl N-(4-chloro-3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}phenyl)carbamate,

a prodrug thereof or a pharmaceutically acceptable salt thereof.

15. (Original) The triazole compound of claim 1, which is

3-chloro-4-[4-methyl-5-(1-phenyl-cyclopropyl)-4H-[1,2,4]triazol-3-yl]-benzamide,

{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzoyl}morpholine hydrochloride,

3-chloro-N-methyl-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide hydrochloride,

3-chloro-N,N-dimethyl-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide hydrochloride,

3-chloro-N-(2-hydroxy-ethyl)-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-N-isopropyl-4-[4-methyl-5-(1-phenylcyclopropyl)4H-[1,2,4]triazol-3-yl]benzamide,

{3-chloro-4-[4-methyl-5-(1-phenyl-cyclopropyl)-4H[1,2,4]triazol-3-yl]benzoyl}piperidine hydrochloride,

{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H[1,2,4]triazol-3-yl]benzoyl}-(4-hydroxy)piperidine,

N-carbamoylmethyl-3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-N-(2,2,2-trifluoro-ethyl)-benzamide hydrochloride,

N-(2-acetylamino)ethyl-3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide hydrochloride,

3-chloro-N-(2-methoxy)ethyl-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide hydrochloride,

1-acetyl-(4-{3-Chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzoyl}piperazine hydrochloride,

3-chloro-N-(2-dimethylamino)ethyl-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-N-(2-morpholin-4-yl)ethylbenzamide,

4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-3-methoxybenzamide,

3-chloro-4-{4-methyl-5-[1-(4-fluorophenyl)cyclopropyl]-4H-[1,2,4]triazol-3-yl}benzamide,

3-chloro-N-metyl-4-{4-methyl-5-[1-(4-fluoro-phenyl)cyclopropyl]-4H-[1,2,4]triazol-3-yl}benzamide,

4-[4-isopropyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-4-[4-methyl-5-(1-thiophen-2-yl)cyclopropyl-4H-[1,2,4]triazol-3-yl]benzamide,

4-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}benzamide,

4-chloro-3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}benzamide,

4-chloro-3-{5-[1-phenylcyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}benzamide,

3-chloro-4-[4-ethyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-chloro-4-{4-ethyl-5-[1-(4-fluorophenyl)cyclopropyl]-4H-[1,2,4]triazol-3-yl}benzamide,

3-[4-isopropyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]benzamide,

3-{5-[1-(4-fluoro-phenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}benzamide,

a prodrug thereof or a pharmaceutically acceptable salt thereof.

16. (Original) The triazole compound of claim 1, which is

N- { 3-chloro-4-[4-methyl-5-(1-phenylcycloproppyl)-4H-[1,2,4]triazol-3-yl]phenyl } -1-morpholinecarboxamide,

3-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}-1,1-dimethylurea,

{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}urea,

ethyl N-{3-Chloro-4-[4-methyl-5-(1-phenyl-cyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}-carbamate,

N- { 3-chloro-4-[4-methyl-5-(1-phenylcycloproppyl)-4H-[1,2,4]triazol-3-yl]phenyl } -1-(4-methoxypiperidine)carboxamide,

N- { 3-chloro-4-[4-methyl-5-(1-phenylcycloproppyl)-4H-[1,2,4]triazol-3-yl]phenyl } -1-(3-hydroxypiperidine)carboxamide,

N- { 3-chloro-4-[4-methyl-5-(1-phenylcycloproppyl)-4H-[1,2,4]triazol-3-yl]phenyl } -1-(4-hydroxypiperidine)carboxamide,

1-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}-3-methoxyurea,

1-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]phenyl}-3-hydroxy-3-methylurea,

1-(3-chloro-4-{5-[1-(4-fluorophenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}phenyl)-3-methoxyurea,

1-(4-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}phenyl)-3-methoxyurea,

1-(3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}phenyl)-3-methoxyurea,

a prodrug thereof or a pharmaceutically acceptable salt thereof.

17. (Original) The triazole compound of claim 1, which is

3-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]-phenyl}oxazolidin-2-one,

1-{3-chloro-4-[4-methyl-5-(1-phenylcyclopropyl)-4H-[1,2,4]triazol-3-yl]phenyl}imidazolidin-2-one,

3-(3-chloro-4-{5-[1-(4-fluoro-phenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}phenyl)oxazolidin-2-one,

3-(4-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}phenyl)oxazolidin-2-one,

3-(4-chloro-3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-methyl-4H-[1,2,4]triazol-3-yl}phenyl)oxazolidin-2-one,

3-(3-{5-[1-(4-fluorophenyl)cyclopropyl]-4-isopropyl-4H-[1,2,4]triazol-3-yl}phenyl)oxazolidin-2-one,

a prodrug thereof or a pharmaceutically acceptable salt thereof.

18. (Previously Presented) A pharmaceutical composition comprising the triazole compound of claim 1, a prodrug thereof or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.

19. (Withdrawn) An HSD1 (11beta-hydroxysteroid dehydrogenase 1) inhibitor comprising the triazole compound of claim 1, a prodrug thereof or a pharmaceutically acceptable salt thereof as an effective component.

20. (Withdrawn) A therapeutic or prophylactic drug of diabetes, which comprises the triazole compound of claim 1, a prodrug thereof or a pharmaceutically acceptable salt thereof as an effective component.

21. (Withdrawn) A therapeutic or prophylactic drug of obesity, which comprises the triazole compound of claim 1, a prodrug thereof or a pharmaceutically acceptable salt thereof as an effective component.

22. (Withdrawn) A therapeutic or prophylactic drug of metabolic syndrome, which comprises the triazole compound of claim 1, a prodrug thereof or a pharmaceutically acceptable salt thereof as an effective component.

23. (Withdrawn) A method for the treatment or prophylaxis of diabetes, which comprises administering an effective amount of the triazole compound of claim 1, a prodrug thereof or a pharmaceutically acceptable salt thereof to a mammal.

24. (Withdrawn) A method for the treatment or prophylaxis of obesity, which comprises administering an effective amount of the triazole compound of claim 1, a prodrug thereof or a pharmaceutically acceptable salt thereof to a mammal.

25. (Withdrawn) A method for the treatment or prophylaxis of metabolic syndrome, which comprises administering an effective amount of the triazole compound of claim 1, a prodrug thereof or a pharmaceutically acceptable salt thereof to a mammal.

26. (Withdrawn) The method of claim 23, wherein a different therapeutic drug of diabetes is used in combination.

27. (Withdrawn) The method of claim 26, wherein the different therapeutic drug of diabetes is one or more pharmaceutical agents selected from the group consisting of an insulin preparation, a sulfonylurea, an insulin secretagogue, a sulfonamide, a biguanide, an α-glucosidase inhibitor and an insulin sensitizer.

28. (Withdrawn) The method of claim 27, wherein the different therapeutic drug of diabetes is one or more pharmaceutical agents selected from the group consisting of insulin, glibenclamide, tolbutamide, glyclopyramide, acetohexamide, glimepiride, tolazamide, gliclazide, nateglinide, glybuzole, metformin hydrochloride, buformine hydrochloride, voglibose, acarbose and pioglitazone hydrochloride.

29. (Withdrawn) The method of claim 24, wherein a different therapeutic drug of diabetes is used in combination.

30. (Withdrawn) The method of claim 29, wherein the different therapeutic drug of diabetes is one or more pharmaceutical agents selected from the group consisting of an insulin preparation, a sulfonylurea, an insulin secretagogue, a sulfonamide, a biguanide, an α-glucosidase inhibitor and an insulin sensitizer.

31. (Withdrawn) The method of claim 30, wherein the different therapeutic drug of diabetes is one or more pharmaceutical agents selected from the group consisting of insulin, glibenclamide, tolbutamide, glyclopyramide, acetohexamide, glimepiride, tolazamide, gliclazide, nateglinide, glybuzole, metformin hydrochloride, buformine hydrochloride, voglibose, acarbose and pioglitazone hydrochloride.

32. (Withdrawn) The method of claim 25, wherein a different therapeutic drug of diabetes is used in combination.

33. (Withdrawn) The method of claim 32, wherein the different therapeutic drug of diabetes is one or more pharmaceutical agents selected from the group consisting of an insulin preparation, a sulfonylurea, an insulin secretagogue, a sulfonamide, a biguanide, an α-glucosidase inhibitor and an insulin sensitizer.

34. (Withdrawn) The method of claim 33, wherein the different therapeutic drug of diabetes is one or more pharmaceutical agents selected from the group consisting of insulin, glibenclamide, tolbutamide, glyclopyramide, acetohexamide, glimepiride, tolazamide, gliclazide, nateglinide, glybuzole, metformin hydrochloride, buformine hydrochloride, voglibose, acarbose and pioglitazone hydrochloride.

35. (Withdrawn) The method of claim 23, wherein a different therapeutic drug of obesity is used in combination.

36. (Withdrawn) The method of claim 35, wherein the different therapeutic drug of obesity is Mazindol.

37. (Withdrawn): The method of claim 24, wherein a different therapeutic drug of obesity is used in combination.

38. (Withdrawn) The method of claim 37, wherein the different therapeutic drug of obesity is Mazindol.

39. (Withdrawn) The method of claim 25, wherein a different therapeutic drug of obesity is used in combination.

40. (Withdrawn) The method of claim 39, wherein the different therapeutic drug of obesity is Mazindol.